

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU090718\
Data File : VU026612.D
Acq On : 07 Sep 2018 10:09
Operator : MD/SY
Sample : J4806-01
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
TAN-BLDG-1FL-KITCHEN-SINK

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U082018DW.M
Title : METHOD 524.2 VOLATILES DRINKING WATER

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.444	384	395	416	rBV	33218	59859	4.22%	3.457%
2	4.675	1070	1089	1116	rBV	622091	1418476	100.00%	81.911%
3	5.746	1408	1422	1437	rBV	27141	53072	3.74%	3.065%
4	6.759	1724	1737	1758	rVB	63784	117381	8.28%	6.778%
5	10.312	2832	2842	2864	rVB2	22108	34647	2.44%	2.001%
6	11.861	3314	3324	3338	rBV2	29766	48296	3.40%	2.789%

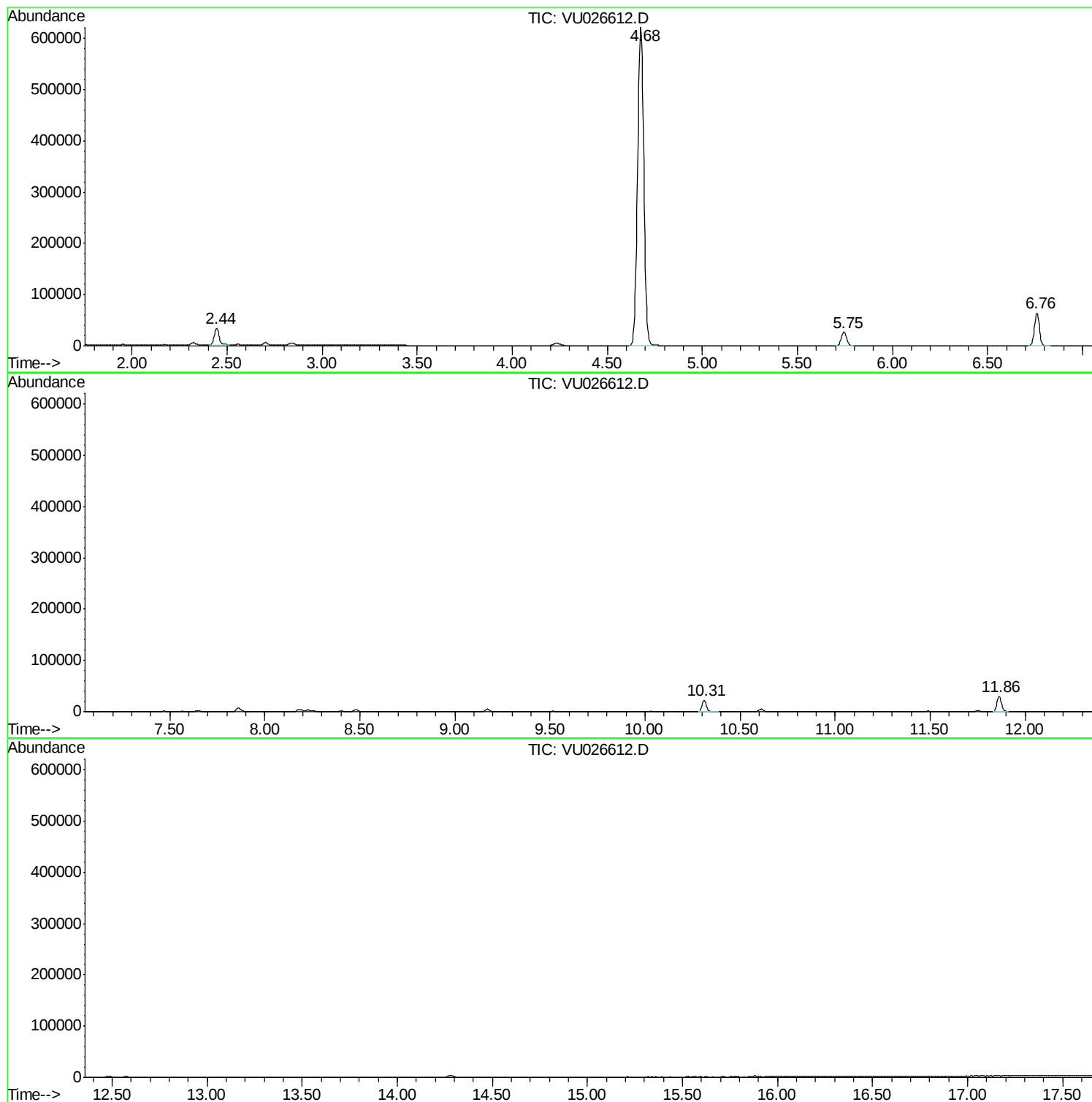
Sum of corrected areas: 1731731

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA U\METHOD\524U082018DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Isopropyl Alcohol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.44	1.13 ug/l	59859	Fluorobenzene	5.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl Alcohol	60	C3H8O	000067-63-0	74
2			Isopropyl Alcohol	60	C3H8O	000067-63-0	74
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	9
4			2-Propanone, 1-methoxy-	88	C4H8O2	005878-19-3	4
5			4-Penten-2-ol	86	C5H10O	000625-31-0	4

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Isopropyl Alcohol	2.44	1.1	ug/l	59859	1	5.75	53072	1.0